

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
 Data File : BP001970.D
 Acq On : 29 Feb 2020 21:53
 Operator : CG/JU
 Sample : L1615-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDP2

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	12	rVB	216745	278833	3.23%	0.321%
2	3.164	43	47	55	rVB	77007	107147	1.24%	0.123%
3	3.899	169	172	176	rVB2	8906	11826	0.14%	0.014%
4	4.793	319	324	332	rBV	132435	189495	2.20%	0.218%
5	6.822	662	669	680	rBV	1277873	2107812	24.43%	2.426%
6	6.981	688	696	711	rVB	1054010	1634761	18.95%	1.882%
7	7.175	722	729	740	rBV	1433767	2265770	26.26%	2.608%
8	7.640	799	808	818	rBV	883835	1381363	16.01%	1.590%
9	7.722	818	822	828	rVB2	9004	14880	0.17%	0.017%
10	8.346	920	928	943	rBV	1636306	2600743	30.15%	2.994%
11	8.781	991	1002	1016	rBV	1177085	1920844	22.27%	2.211%
12	9.275	1081	1086	1093	rVB	34311	53637	0.62%	0.062%
13	9.499	1117	1124	1142	rBV	947314	1590210	18.43%	1.830%
14	9.928	1189	1197	1201	rBV8	4928	11147	0.13%	0.013%
15	10.040	1208	1216	1233	rBV	1667180	2841816	32.94%	3.271%
16	10.410	1271	1279	1290	rBV	1156518	1964393	22.77%	2.261%
17	10.546	1294	1302	1322	rBV	1417480	2476317	28.70%	2.850%
18	11.951	1533	1541	1547	rBV	79853	151216	1.75%	0.174%
19	12.010	1547	1551	1557	rVB	26755	46666	0.54%	0.054%
20	13.022	1716	1723	1728	rBV4	31210	51482	0.60%	0.059%
21	13.134	1737	1742	1744	rBV2	10563	13198	0.15%	0.015%
22	13.175	1744	1749	1757	rVB6	28534	62012	0.72%	0.071%
23	13.693	1830	1837	1842	rBV	2876241	3989512	46.24%	4.592%
24	13.734	1842	1844	1853	rVB	220717	295219	3.42%	0.340%
25	13.963	1876	1883	1897	rBV	3361139	5077267	58.85%	5.844%
26	14.193	1917	1922	1930	rVB2	18305	31621	0.37%	0.036%
27	14.275	1930	1936	1947	rBV	1845140	2735893	31.71%	3.149%
28	14.369	1947	1952	1961	rVB3	19282	31498	0.37%	0.036%
29	14.475	1964	1970	1991	rBV	981061	1584677	18.37%	1.824%
30	14.704	2004	2009	2012	rBV7	5306	9131	0.11%	0.011%
31	15.175	2083	2089	2093	rBV2	12367	19290	0.22%	0.022%
32	15.275	2097	2106	2113	rBV	4572500	6356607	73.68%	7.317%
33	15.392	2120	2126	2142	rBV	929982	1330523	15.42%	1.531%
34	15.516	2142	2147	2153	rVB	50591	74266	0.86%	0.085%

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35	15.769	2186	2190	2196	rVB9	7085	14127	0.16%	0.016%
36	15.975	2219	2225	2233	rBV	126058	202265	2.34%	0.233%
37	16.081	2237	2243	2249	rVB3	34513	71631	0.83%	0.082%
38	16.292	2276	2279	2284	rVB	8665	11256	0.13%	0.013%
39	16.710	2346	2350	2354	rVB3	10626	15291	0.18%	0.018%
40	16.822	2363	2369	2377	rBV2	10360	23804	0.28%	0.027%
41	17.028	2390	2404	2414	rBV	2308039	3319623	38.48%	3.821%
42	17.128	2414	2421	2433	rBV	4707720	6663623	77.24%	7.670%
43	17.398	2464	2467	2471	rVV5	4592	9271	0.11%	0.011%
44	17.445	2471	2475	2484	rVB3	22465	39254	0.46%	0.045%
45	17.716	2516	2521	2528	rVB3	22333	45478	0.53%	0.052%
46	17.904	2549	2553	2556	rBV6	8547	12598	0.15%	0.015%
47	17.951	2556	2561	2568	rBV	226365	345110	4.00%	0.397%
48	18.098	2582	2586	2593	rVB3	58454	89062	1.03%	0.103%
49	18.239	2606	2610	2615	rBV2	26415	39879	0.46%	0.046%
50	18.375	2629	2633	2638	rBV6	13304	17193	0.20%	0.020%
51	18.786	2701	2703	2712	rVB10	16997	26664	0.31%	0.031%
52	18.945	2728	2730	2734	rVB5	6953	10198	0.12%	0.012%
53	19.016	2738	2742	2746	rBV	61440	89010	1.03%	0.102%
54	19.051	2746	2748	2750	rVV2	16610	18669	0.22%	0.021%
55	19.075	2750	2752	2759	rVB8	13833	23897	0.28%	0.028%
56	19.192	2768	2772	2780	rBV	61470	102606	1.19%	0.118%
57	19.263	2780	2784	2790	rVB	61608	83225	0.96%	0.096%
58	19.375	2797	2803	2809	rBV2	5924626	8077908	93.63%	9.298%
59	19.451	2814	2816	2823	rVB6	25553	27091	0.31%	0.031%
60	19.616	2842	2844	2854	rVB9	27783	40442	0.47%	0.047%
61	19.939	2896	2899	2903	rVB	27203	28985	0.34%	0.033%
62	20.422	2978	2981	2984	rVB	59387	60328	0.70%	0.069%
63	20.869	3052	3057	3062	rBV2	493770	771464	8.94%	0.888%
64	21.122	3095	3100	3109	rBV	3198883	3933130	45.59%	4.527%
65	21.239	3117	3120	3126	rBV	58572	99177	1.15%	0.114%
66	21.298	3127	3130	3135	rVV	283067	329418	3.82%	0.379%
67	21.733	3200	3204	3215	rVB	479801	602603	6.98%	0.694%
68	22.192	3278	3282	3290	rVB	634543	825439	9.57%	0.950%
69	22.322	3299	3304	3311	rBV	257616	356011	4.13%	0.410%
70	22.698	3363	3368	3378	rVB	732344	1121821	13.00%	1.291%
71	23.192	3444	3452	3460	rBV	4713805	8627136	100.00%	9.930%

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Title : SVOA CALIBRATION

72	23.269	3460	3465	3469	rVV	631967	1093873	12.68%	1.259%
73	23.333	3469	3476	3489	rVB	2189939	4178524	48.43%	4.810%
74	23.916	3569	3575	3583	rBV	505949	986063	11.43%	1.135%
75	24.668	3697	3703	3719	rVB2	304585	761997	8.83%	0.877%
76	25.545	3846	3852	3866	rVB2	131349	370907	4.30%	0.427%

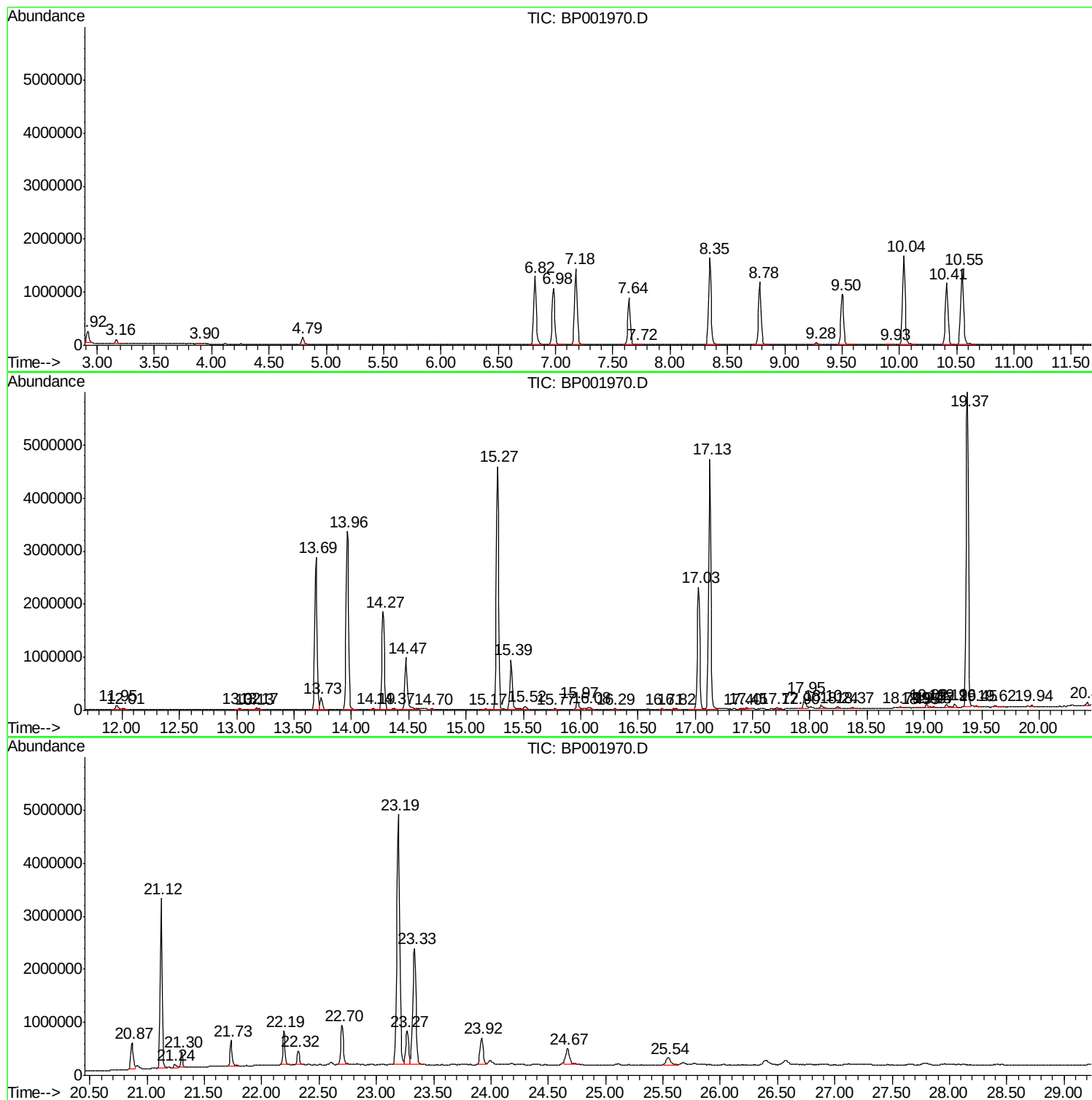
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Instrument :
BNA_P
ClientSampled :
DBDP2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022920\
Data File : BP001970.D
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Instrument :
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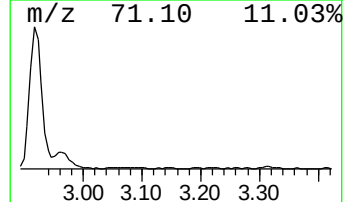
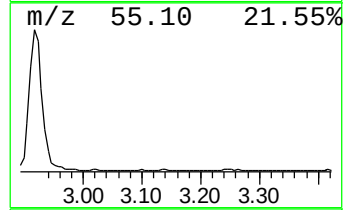
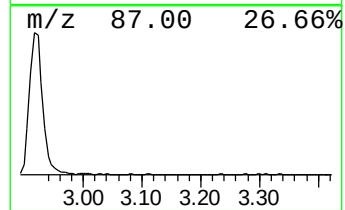
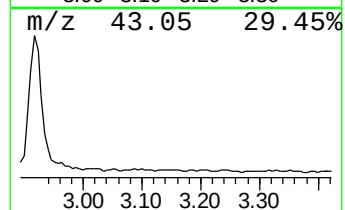
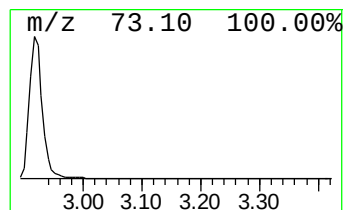
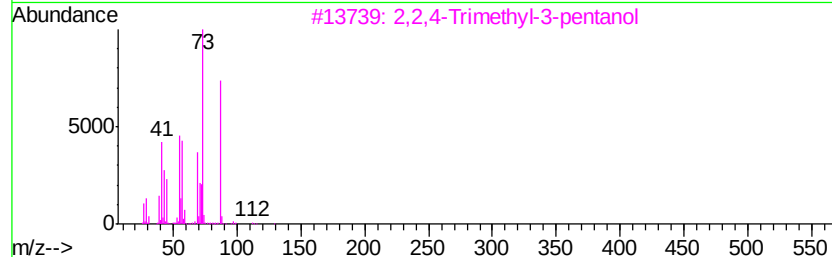
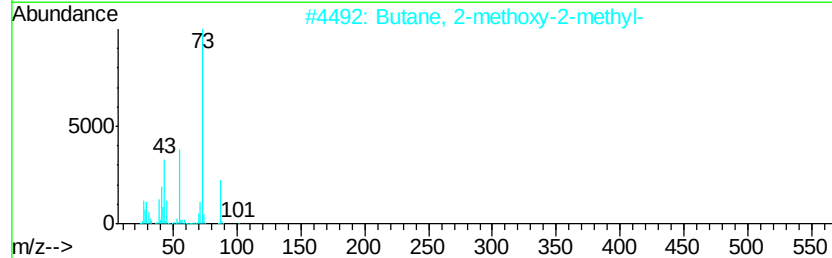
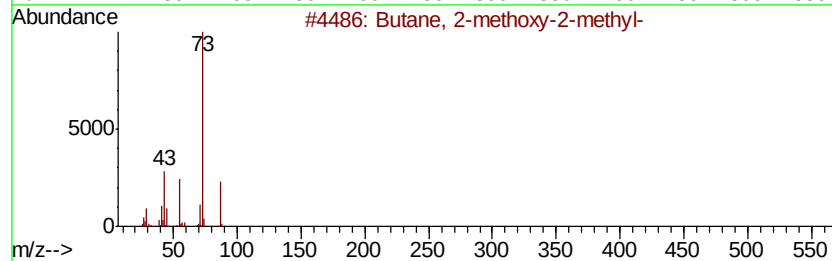
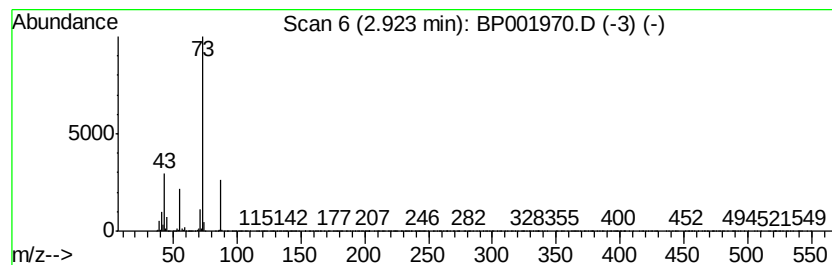
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	4.04 ng/ul	278833	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	16



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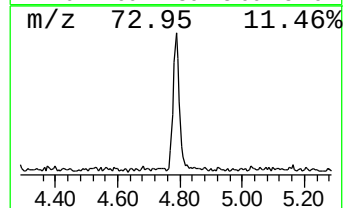
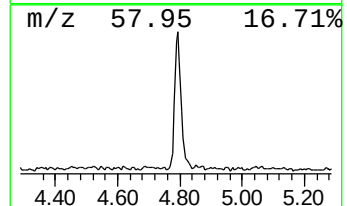
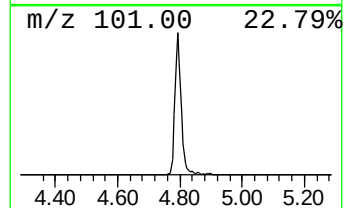
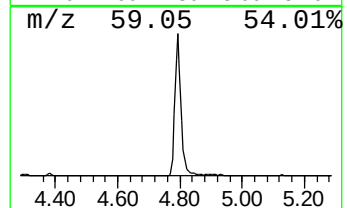
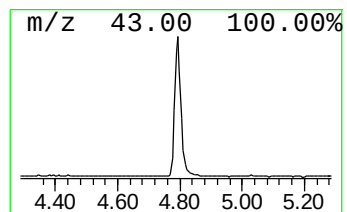
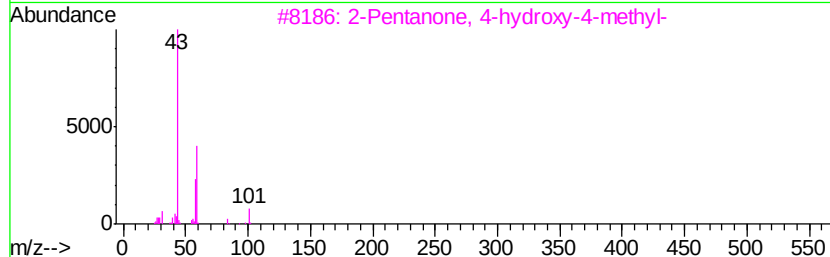
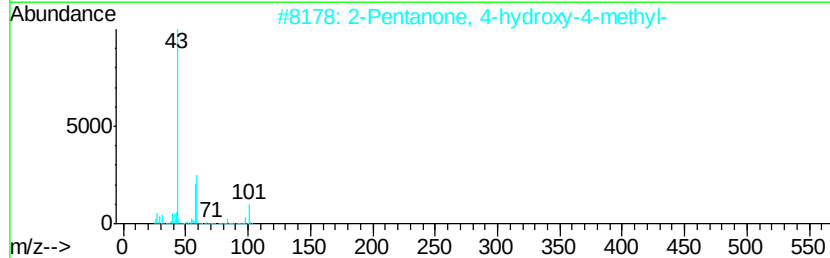
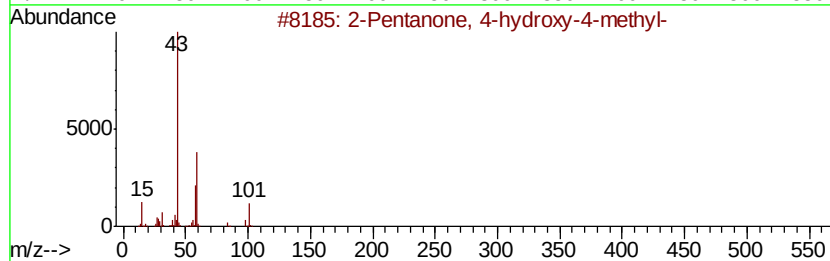
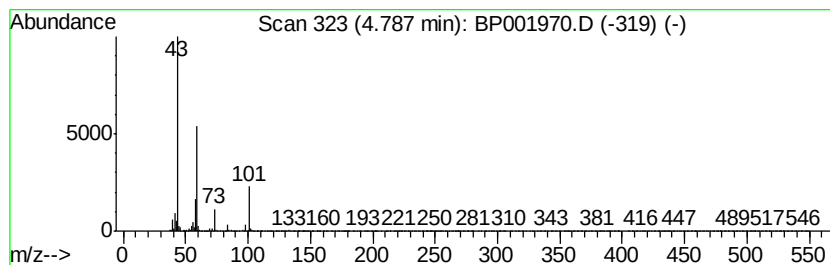
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	2.74 ng/ul	189495	1,4-Dichlorobenzene-d4	7.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4		2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	25
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	25



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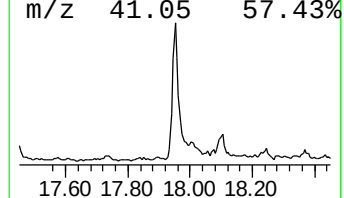
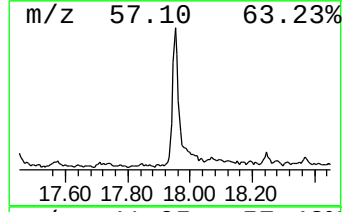
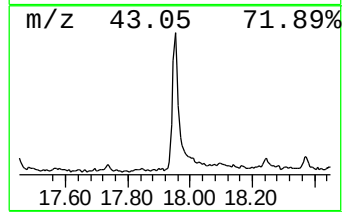
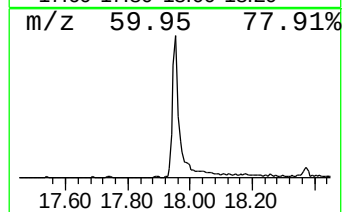
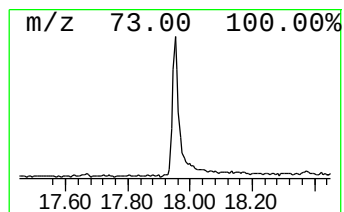
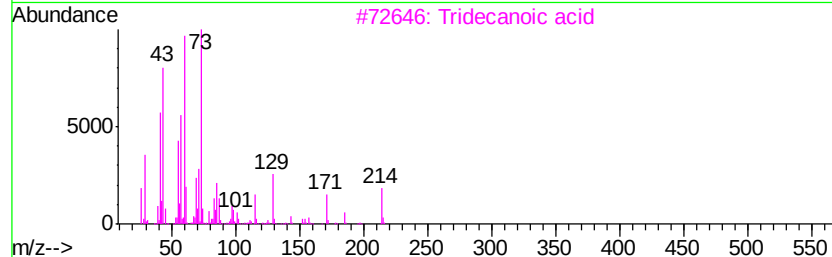
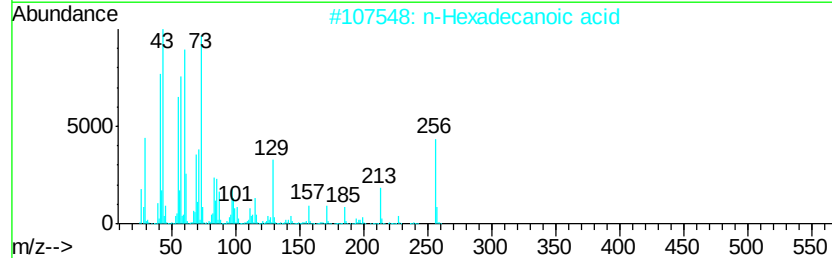
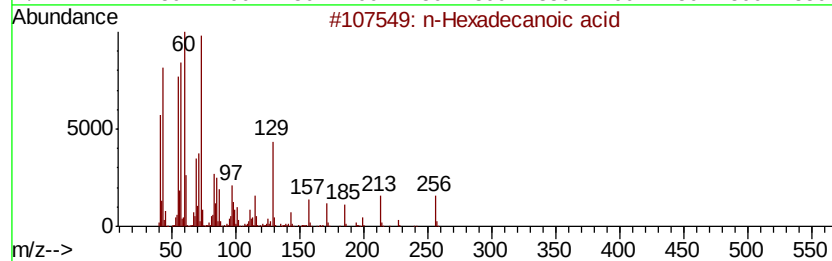
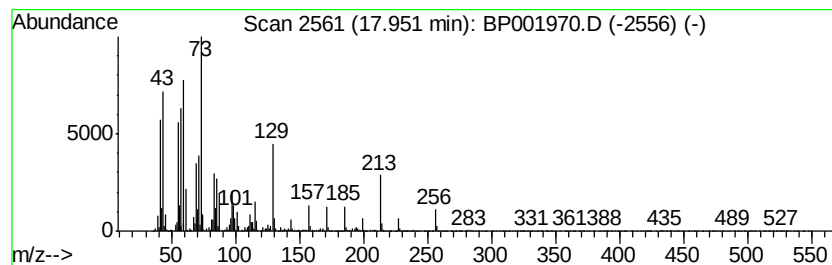
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 3 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
17.95	2.08 ng/ul	345110	Phenanthrene-d10		17.03		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			Tridecanoic acid	214	C13H26O2	000638-53-9	91
4			Octadecanoic acid	284	C18H36O2	000057-11-4	87
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	86



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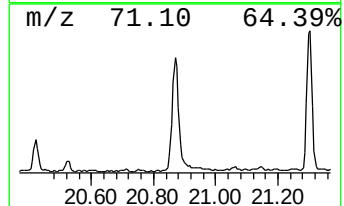
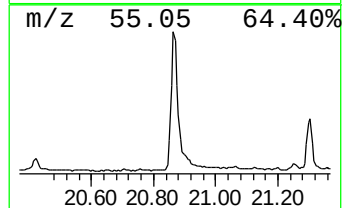
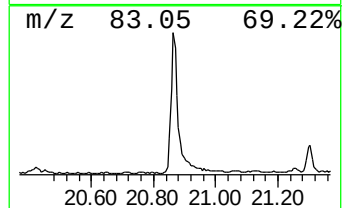
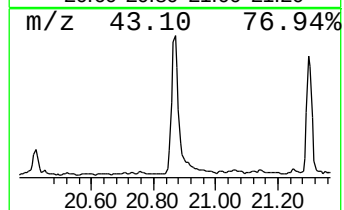
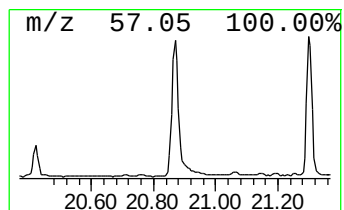
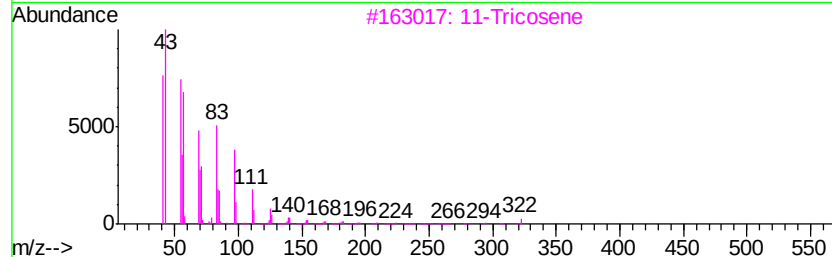
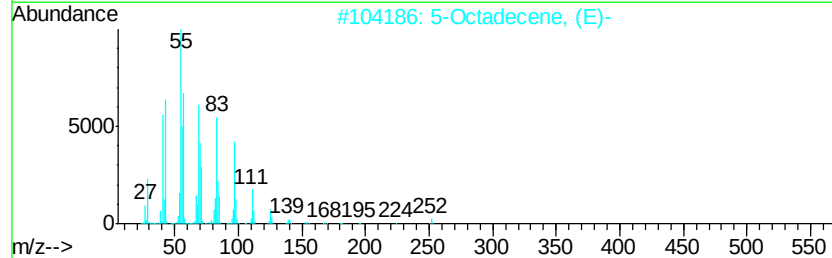
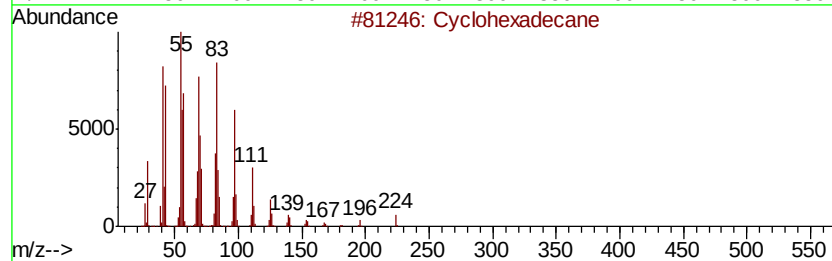
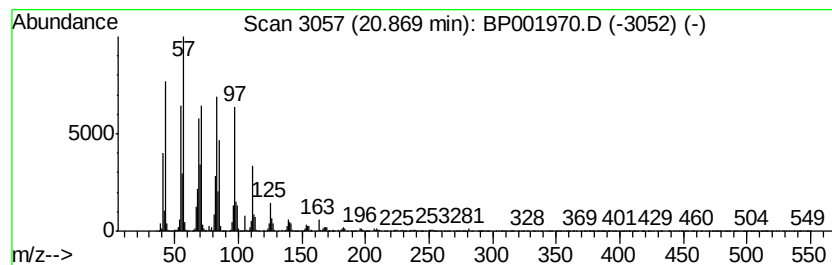
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Cyclic20.87 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	3.92 ng/ul	771464	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	99
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	95
3		11-Tricosene	322	C23H46	052078-56-5	92
4		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5		Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
Data File : BP001970.D
Acq On : 29 Feb 2020 21:53
Operator : CG/JU
Sample : L1615-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBDP2

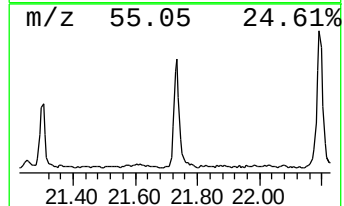
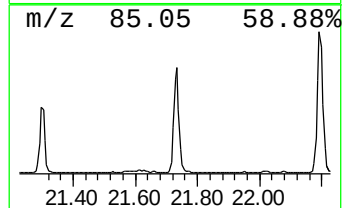
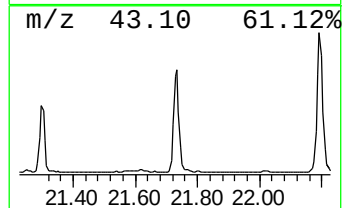
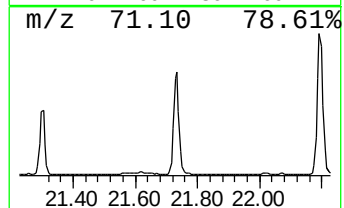
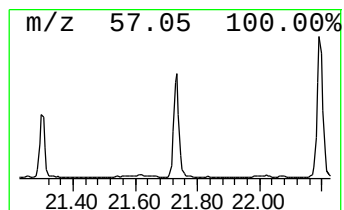
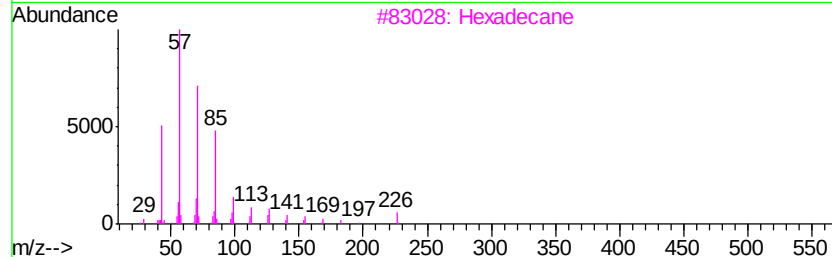
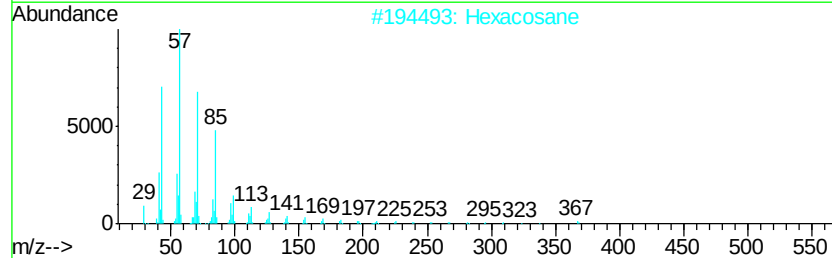
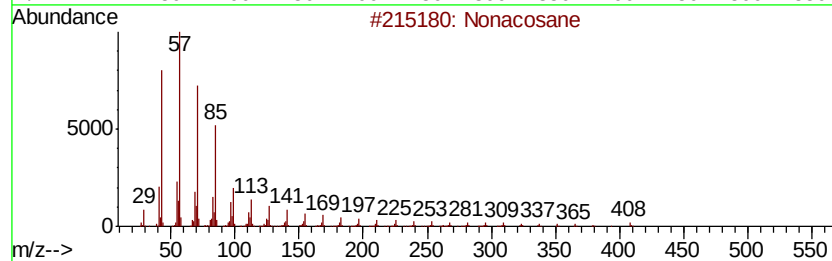
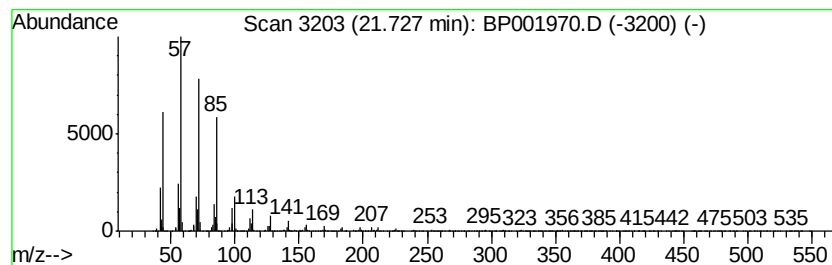
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.73	3.06 ng/ul	602603	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonacosane	408	C29H60	000630-03-5	97
2		Hexacosane	366	C26H54	000630-01-3	94
3		Hexadecane	226	C16H34	000544-76-3	93
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
Data File : BP001970.D
Acq On : 29 Feb 2020 21:53
Operator : CG/JU
Sample : L1615-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBDP2

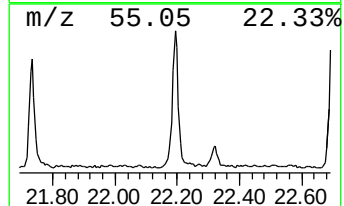
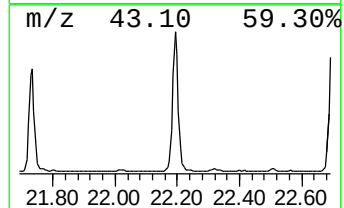
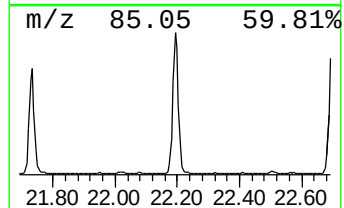
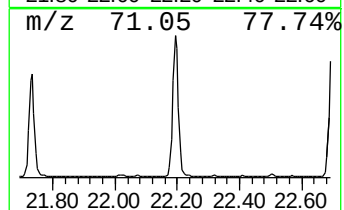
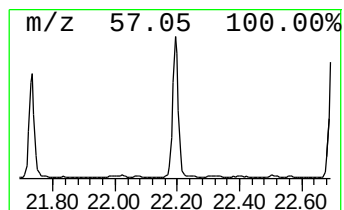
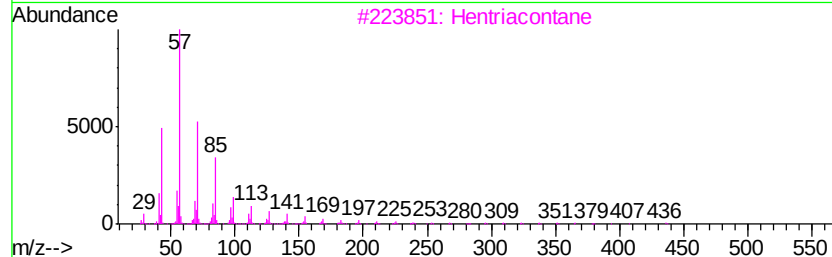
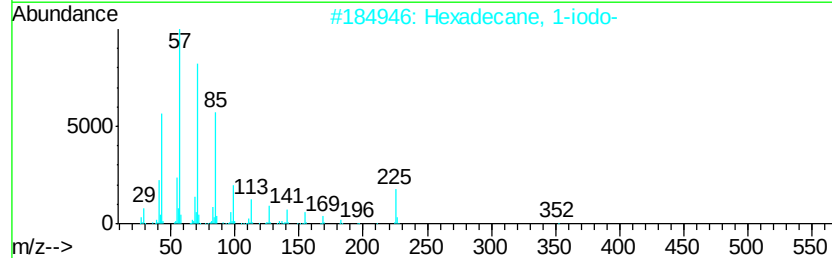
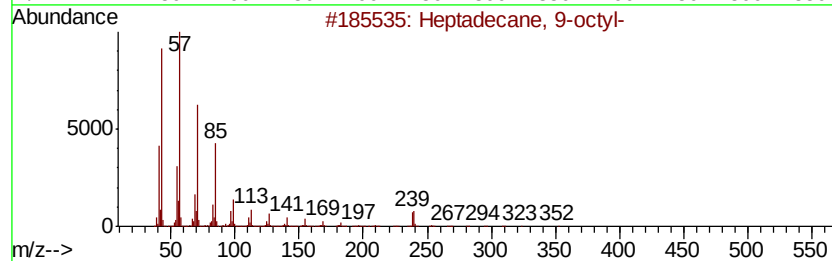
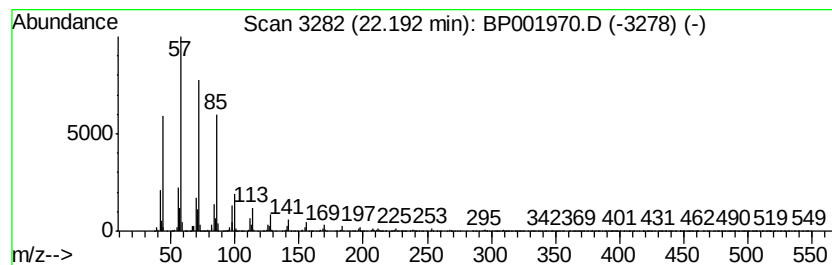
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.19	4.20 ng/ul	825439	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	97
2		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
3		Hentriacontane	437	C31H64	000630-04-6	93
4		Tetratriacontane	479	C34H70	014167-59-0	93
5		Pentacosane	352	C25H52	000629-99-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
Data File : BP001970.D
Acq On : 29 Feb 2020 21:53
Operator : CG/JU
Sample : L1615-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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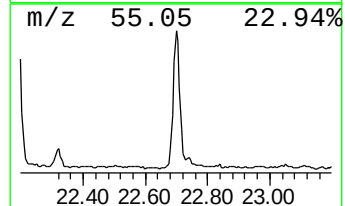
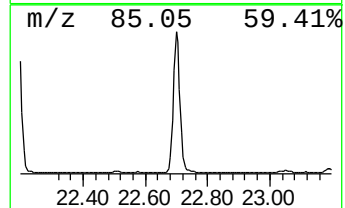
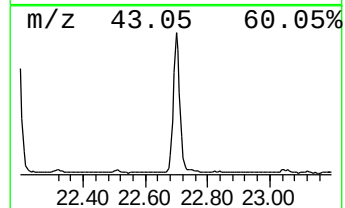
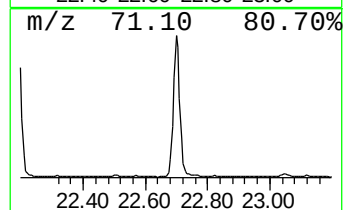
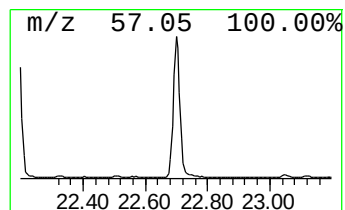
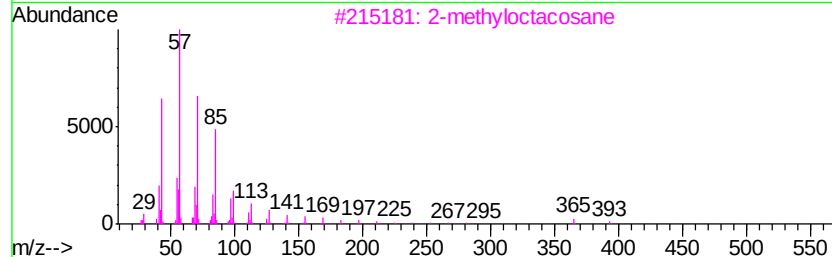
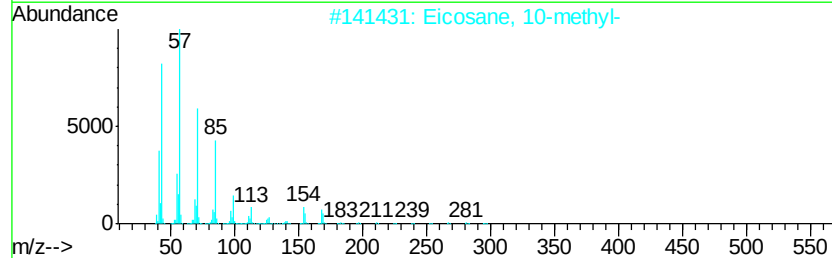
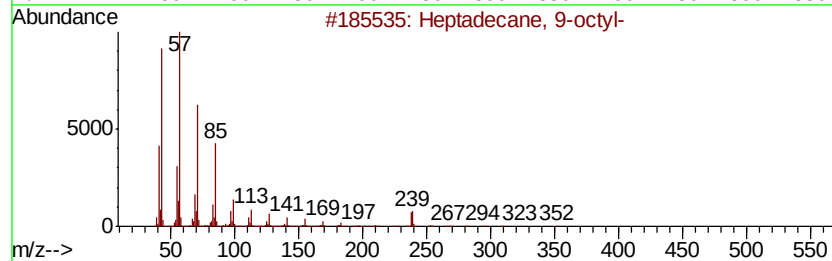
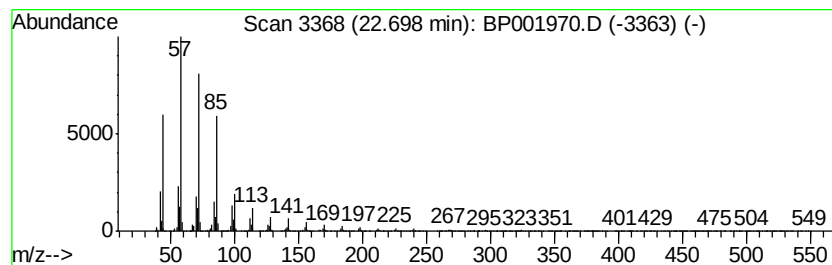
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.70	5.37 ng/ul	1121820	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
Data File : BP001970.D
Acq On : 29 Feb 2020 21:53
Operator : CG/JU
Sample : L1615-18
Misc :
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Instrument :
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ClientSampleId :
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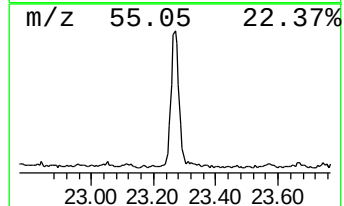
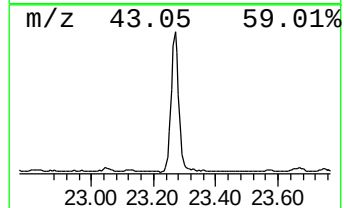
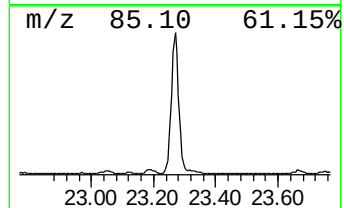
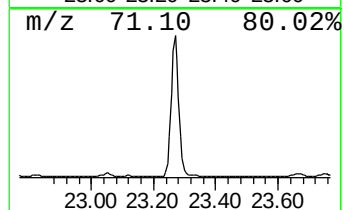
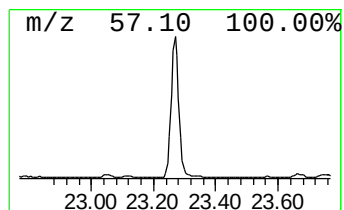
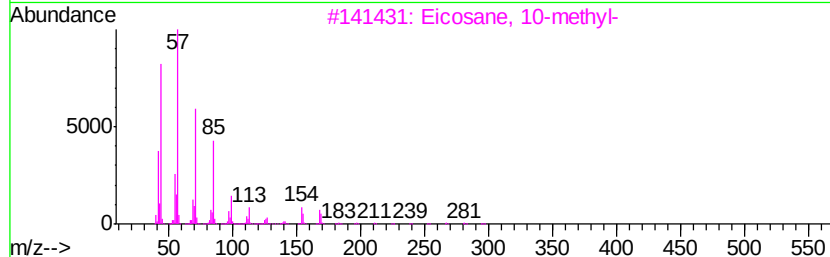
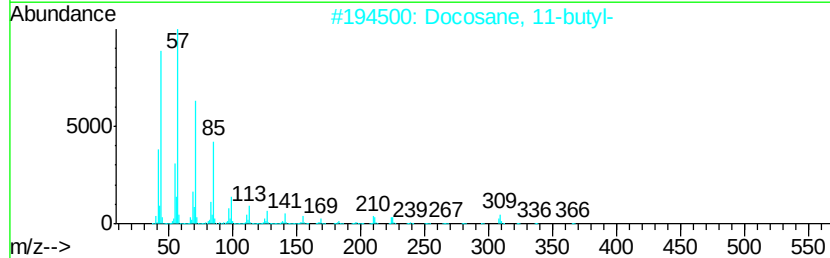
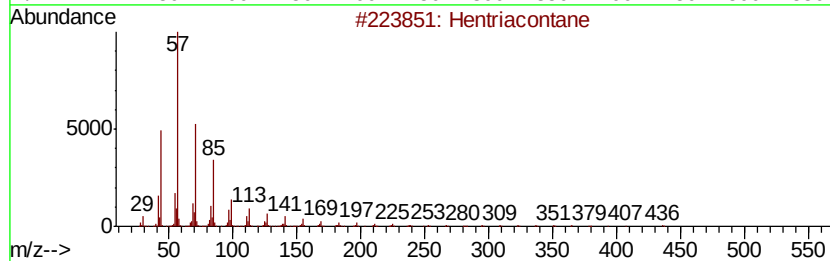
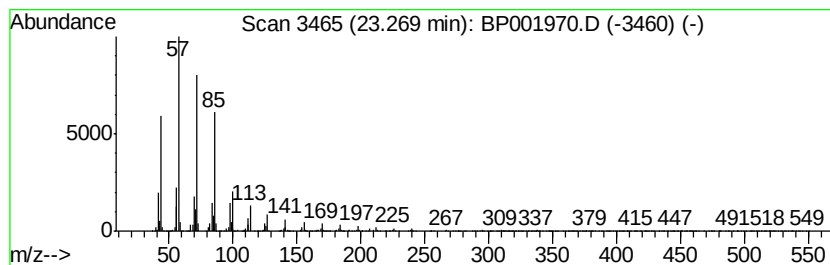
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.27	5.24 ng/ul	1093870	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	96
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
Data File : BP001970.D
Acq On : 29 Feb 2020 21:53
Operator : CG/JU
Sample : L1615-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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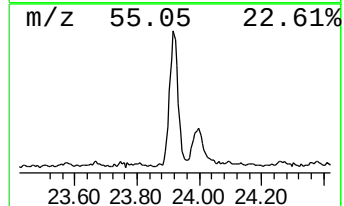
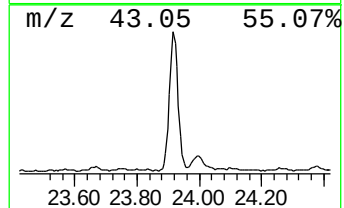
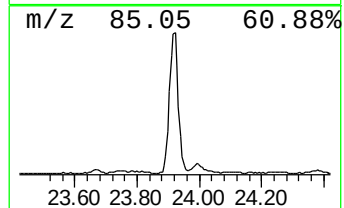
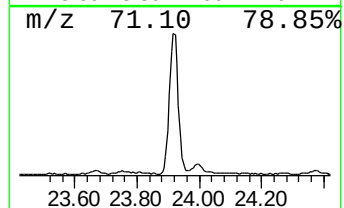
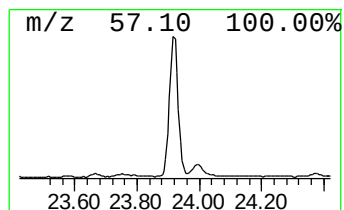
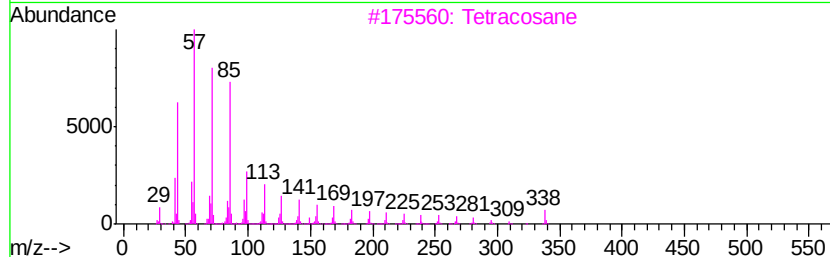
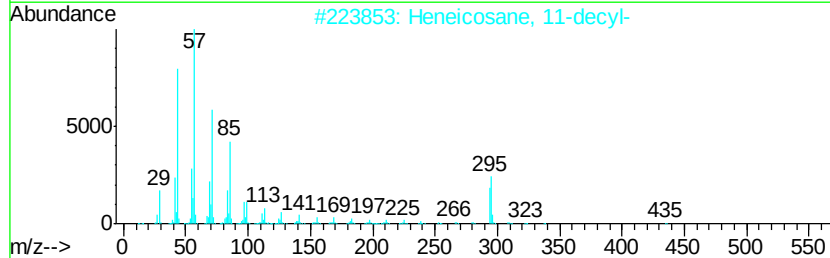
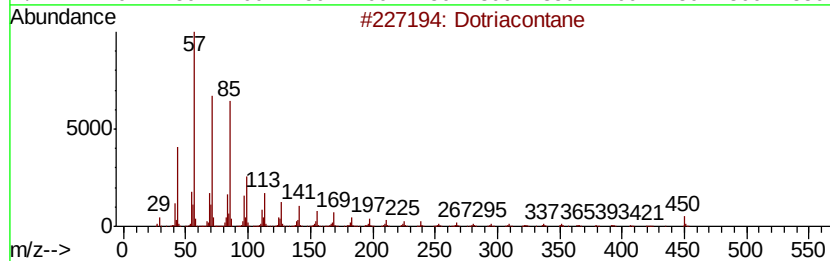
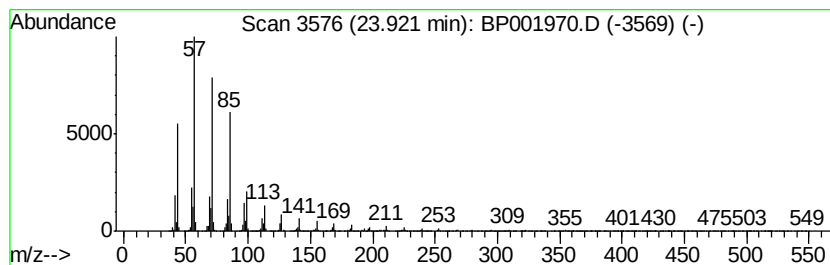
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.92	4.72 ng/ul	986063	Perylene-d12	23.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dotriacontane	451	C32H66	000544-85-4	97
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95
3		Tetracosane	338	C24H50	000646-31-1	94
4		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022920\
Data File : BP001970.D
Acq On : 29 Feb 2020 21:53
Operator : CG/JU
Sample : L1615-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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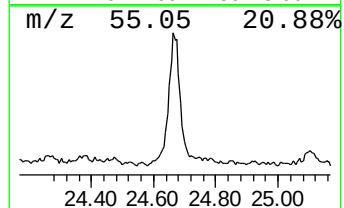
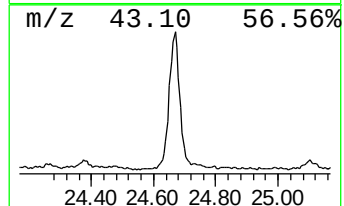
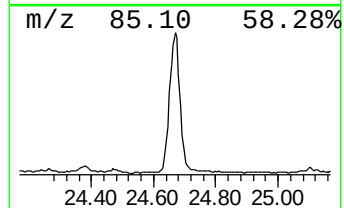
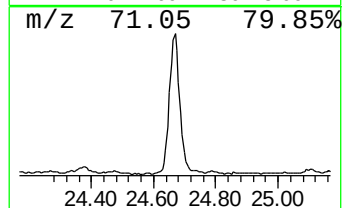
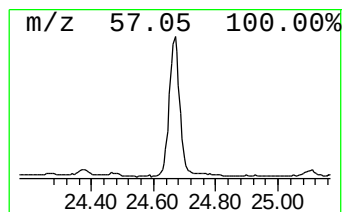
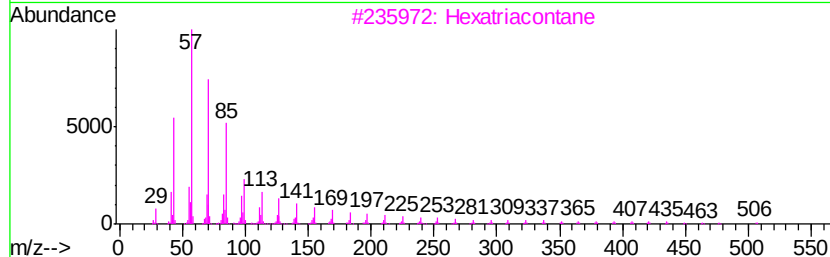
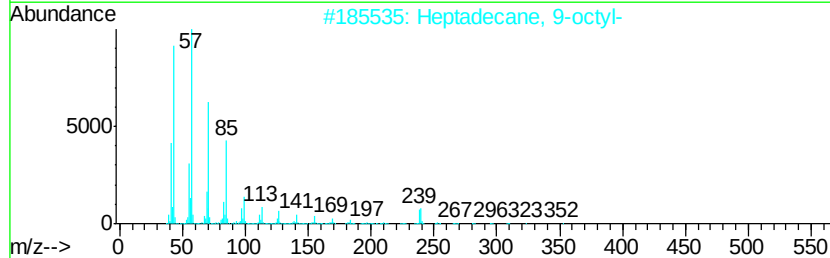
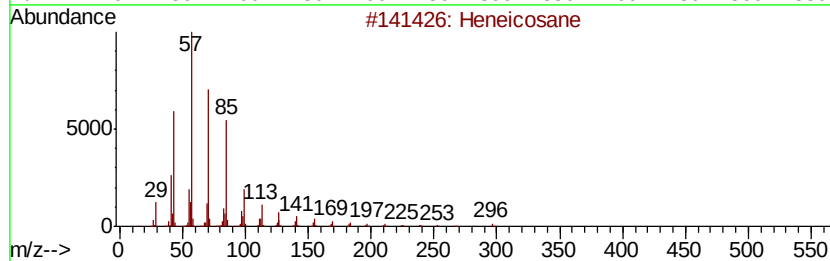
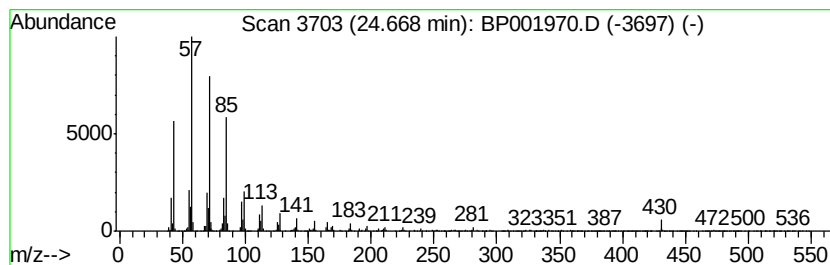
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.67	3.65 ng/ul	761997	Perylene-d12	23.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3			Hexatriacontane	507	C36H74	000630-06-8	93
4			Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
5			Docosane, 7-hexyl-	394	C28H58	055373-86-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022920\
 Data File : BP001970.D
 Acq On : 29 Feb 2020 21:53
 Operator : CG/JU
 Sample : L1615-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDP2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.92	4.0	ng/ul	278833	1	7.64	1381360	20.0
2-Pentanone, 4-hy...	4.79	2.7	ng/ul	189495	1	7.64	1381360	20.0
n-Hexadecanoic acid	17.95	2.1	ng/ul	345110	4	17.03	3319620	20.0
(DEL) Alkane: Cyc...	20.87	3.9	ng/ul	771464	5	21.12	3933130	20.0
(DEL) Alkane: Str...	21.73	3.1	ng/ul	602603	5	21.12	3933130	20.0
(DEL) Alkane: Str...	22.19	4.2	ng/ul	825439	5	21.12	3933130	20.0
(DEL) Alkane: Str...	22.70	5.4	ng/ul	1121820	6	23.33	4178520	20.0
(DEL) Alkane: Str...	23.27	5.2	ng/ul	1093870	6	23.33	4178520	20.0
(DEL) Alkane: Str...	23.92	4.7	ng/ul	986063	6	23.33	4178520	20.0
(DEL) Alkane: Str...	24.67	3.6	ng/ul	761997	6	23.33	4178520	20.0